

## 90358

Access DB#\_\_\_\_\_

### SEARCH REQUEST FORM 4

#### Scientific and Technical Information Center

Requester's Full Name: BEN SACKEY Examiner #: 73  Art Unit: 16 20 Phone Number 30 5 - 6889 Serial Numb  Mail Box and Bldg/Room Location: CM 3819 Results Format Preferre	489 Date: 3/3/03  per: 10 / 049, 463  ped (circle): PAPER DISK E-MAIL
If more than one search is submitted, please prioritize searches in ord	
Please provide a detailed statement of the search topic, and describe as specifically as possi Include the elected species or structures, keywords, synonyms, acronyms, and registry numutility of the invention. Define any terms that may have a special meaning. Give examples known. Please attach a copy of the cover sheet, pertinent claims, and abstract.	bers, and combine with the concept or
Title of Invention: Method for making methyl ethyl	Ketone Cyanolydvin
Inventors (please provide full names): (Noi 27 et d	
Earliest Priority Filing Date: 7 / 29 / 99	
*For Sequence Searches Only* Please include all pertinent information (parent, child, divisional appropriate serial number.  X yand a cid + methyl ethyl ketone che	•
Jan Delavai Reference Librarian Biotechnology & Chemical Library CM1 1E07 – 703-308-4498 jan.delaval@uspto.gov	TECEIVED  (SAR 31 ZE:  (STIC)
	•

STAFF USE ONLY	Type of Search	Vendors and cost where applicable	
Searcher:	NA Sequence (#)	STN	
Searcher Phone #: 4496	AA Sequence (#)	Dialog	
Searcher Location:	Structure (#)	Questel/Orbit	
Date Searcher Picked Up: 47 03	Bibliographic	Dr.Link	
Date Completed: 412 03	Litigation	Lexis/Nexis	
	Fulltext	Sequence Systems	
Clerical Prep Time:	Patent Family	WWW/Internet	
Online Time:	Other	Other (specify)	
PTO-1590 (8-01)			

# BioTech-Chem Library Search Results Feedback Form (Optional)



The search results generated for your recent request are attached. If you have any questions or comments (compliments or complaints) about the scope or the results of the search, please contact *the BioTech-Chem searcher* who conducted the search *or contact*:

Mary Hale, Supervisor, 308-4258 CM-1 Room 1E01

'olu	ntary Results Feedback Form
>	I am an examiner in Workgroup: (Example: 1610)
>	Relevant prior art found, search results used as follows:
	102 rejection
	103 rejection
	Cited as being of interest.
	Helped examiner better understand the invention.
	Helped examiner better understand the state of the art in their technology.
	Types of relevant prior art found:
	Foreign Patent(s)
	Non-Patent Literature (journal articles, conference proceedings, new product announcements etc.)
>	Relevant prior art <b>not found</b> :
	Results verified the lack of relevant prior art (helped determine patentability).
	Search results were not useful in determining patentability or understanding the invention.
Otho	er Comments:
	,
	65 Lead forms at the Circulation Desk CM-1 or send to Mary Hale, CM1-1E01 or e-mail

Drop off completed forms at the Circulation Desk CM-1, or send to Mary Hale, CM1-1E01 or e-mail mary.hale@uspto.gov.

=> d his

```
(FILE 'REGISTRY' ENTERED AT 17:17:38 ON 02 APR 2003)
                                                                          Jan Dolaval
                 DEL HIS
                                                                       Reference Librarian
                 E C5H9NO/MF
                                                                  Biotechnology & Chambel 1 (2004)
CM1 1807 - Tausto 1705
             469 S E3
L1
L2
             242 S L1 AND NR>=1
                                                                      jende die Garpage /
L3
             227 S L1 NOT L2
              30 S L3 AND ?CYAN?/CNS
T.4
L5
              48 S L3 AND ?NITRIL?/CNS
L6
              68 S L4, L5
L7
               3 S L6 AND BUTANENITRILE AND 2 HYDROXY 2 METHYL
                 E METHYL ETHYL KETONE/CN
               1 S E3
1.8
                 E DIETHYLAMINE/CN
               1 S E3
1.9
                 E HYDROCYANIC ACID/CN
               1 S E3
L10
                 SEL RN L7
               1 S E1-E3/CRN
L11
     FILE 'HCAPLUS' ENTERED AT 17:20:18 ON 02 APR 2003
L12
              60 S L7
L13
              11 S L7/P
L14
          24426 S L8 OR METHYLETHYLKETONE OR METHYL ETHYLKETONE OR METHYLETHYL
          27282 S L10 OR HCN OR HYDROCYANIC ACID OR HYDROGEN CYANIDE
L15
          20861 S L9 OR DIETHYLAMINE OR (DIETHYL OR DI ETHYL OR DIET) () AMINE
L16
L17
              25 S L12 AND L14-L16
L18
              10 S L13 AND L17
L19
              10 S L13 AND L14
L20
               7 S L13 AND L15
L21
              1 S L13 AND L16
L22
              1 S L21 AND L19, L20
L23
              10 S L13, L18-L21 NOT L22
=> fil reg
```

FILE 'REGISTRY' ENTERED AT 17:23:44 ON 02 APR 2003 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2003 American Chemical Society (ACS)

Property values tagged with IC are from the  ${\tt ZIC/VINITI}$  data file provided by  ${\tt InfoChem.}$ 

STRUCTURE FILE UPDATES: 1 APR 2003 HIGHEST RN 501325-53-7 DICTIONARY FILE UPDATES: 1 APR 2003 HIGHEST RN 501325-53-7

TSCA INFORMATION NOW CURRENT THROUGH MAY 20, 2002

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. See HELP PROPERTIES for more information. See STNote 27, Searching Properties in the CAS Registry File, for complete details: http://www.cas.org/ONLINE/STN/STNOTES/stnotes27.pdf

=> d ide can tot 17

L7 ANSWER 1 OF 3 REGISTRY COPYRIGHT 2003 ACS RN 174849-22-0 REGISTRY

CN Butanenitrile, 2-hydroxy-2-methyl-, (2S)- (9CI) (CA INDEX NAME) OTHER CA INDEX NAMES:

CN Butanenitrile, 2-hydroxy-2-methyl-, (S)-

FS STEREOSEARCH

MF C5 H9 N O

SR CA

LC STN Files: CA, CAPLUS

Absolute stereochemistry.



#### \*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

2 REFERENCES IN FILE CA (1962 TO DATE)
2 REFERENCES IN FILE CAPLUS (1962 TO DATE)

REFERENCE 1: 134:109910

REFERENCE 2: 124:224570

L7 ANSWER 2 OF 3 REGISTRY COPYRIGHT 2003 ACS

RN 122045-29-8 REGISTRY

CN Butanenitrile, 2-hydroxy-2-methyl-, (2R)- (9CI) (CA INDEX NAME)

OTHER CA INDEX NAMES:

CN Butanenitrile, 2-hydroxy-2-methyl-, (R)-

OTHER NAMES:

CN (R)-Butan-2-one cyanohydrin

FS STEREOSEARCH

MF C5 H9 N O

SR CA

LC STN Files: AGRICOLA, BEILSTEIN\*, CA, CAPLUS, CASREACT, GMELIN\* (\*File contains numerically searchable property data)

Absolute stereochemistry.



#### \*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

5 REFERENCES IN FILE CA (1962 TO DATE)

5 REFERENCES IN FILE CAPLUS (1962 TO DATE)

REFERENCE 1: 134:109910

REFERENCE 2: 119:43947

REFERENCE 3: 118:58234

REFERENCE 4: 115:91338

REFERENCE 5: 111:173178

```
sackey - 10 / 049463
L7
     ANSWER 3 OF 3 REGISTRY COPYRIGHT 2003 ACS
RN
     4111-08-4 REGISTRY
CN
     Butanenitrile, 2-hydroxy-2-methyl- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
    Butyronitrile, 2-hydroxy-2-methyl- (6CI, 7CI, 8CI)
OTHER NAMES:
CN
     2-Butanone, cyanohydrin
CN
     2-Hydroxy-2-methylbutanenitrile
CN
     2-Hydroxy-2-methylbutyronitrile
CN
     2-Methyl-2-hydroxybutyronitrile
CN
     Butanone cyanohydrin
CN
     Ethyl methyl ketone cyanohydrin
CN
     Methyl ethyl ketone cyanohydrin
FS
     3D CONCORD
DR
     73683-34-8
MF
     C5 H9 N O
CI
     COM
       N Files: BEILSTEIN*, BIOSIS, CA, CAOLD, CAPLUS, CASREACT, CHEMCATS, CHEMLIST, GMELIN*, IFICDB, IFIPAT, IFIUDB, MEDLINE, TOXCENTER, USPATFULL
LC
     STN Files:
          (*File contains numerically searchable property data)
     Other Sources: EINECS**, NDSL**, TSCA**
          (**Enter CHEMLIST File for up-to-date regulatory information)
   OH
Me-C-Et
   CN
**PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT**
               52 REFERENCES IN FILE CA (1962 TO DATE)
```

```
52 REFERENCES IN FILE CAPLUS (1962 TO DATE)
               7 REFERENCES IN FILE CAOLD (PRIOR TO 1967)
REFERENCE
            1: 137:321544
REFERENCE
            2: 136:247644
REFERENCE
            3:
               135:107069
REFERENCE
               135:107068
            4:
REFERENCE
            5:
               135:29904
REFERENCE
            6:
               134:233091
REFERENCE
           7:
               134:131253
REFERENCE
           8: 134:109910
REFERENCE
           9: 133:146265
REFERENCE 10: 130:324940
```

=> d ide can 18

L8 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2003 ACS RN 78-93-3 REGISTRY

```
sackey - 10 / 049463
CN
     2-Butanone (8CI, 9CI) (CA INDEX NAME)
OTHER NAMES:
CN
     3-Butanone
CN
     Butanone
CN
     Ethyl methyl ketone
CN
CN
     Methyl ethyl ketone
FS
     3D CONCORD
DR
     135311-02-3
MF
     C4 H8 O
CI
     COM
LC
                AGRICOLA, ANABSTR, AQUIRE, BEILSTEIN*, BIOBUSINESS, BIOSIS,
     STN Files:
       BIOTECHNO, CA, CANCERLIT, CAOLD, CAPLUS, CASREACT, CBNB, CEN, CHEMCATS,
       CHEMINFORMRX, CHEMLIST, CHEMSAFE, CIN, CSCHEM, CSNB, DDFU, DETHERM*,
       DIPPR*, DRUGU, EMBASE, ENCOMPLIT, ENCOMPLIT2, ENCOMPPAT2,
      GMELIN*, HODOC*, HSDB*, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE, MRCK*,
       MSDS-OHS, NAPRALERT, NIOSHTIC, PDLCOM*, PIRA, PROMT, RTECS*, SPECINFO,
       SYNTHLINE, TOXCENTER, TULSA, ULIDAT, USPAT2, USPATFULL, VTB
         (*File contains numerically searchable property data)
     Other Sources: DSL**, EINECS**, TSCA**
         (**Enter CHEMLIST File for up-to-date regulatory information)
H3C-C-CH2-CH3
**PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT**
           19505 REFERENCES IN FILE CA (1962 TO DATE)
            173 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
           19529 REFERENCES IN FILE CAPLUS (1962 TO DATE)
             10 REFERENCES IN FILE CAOLD (PRIOR TO 1967)
```

REFERENCE 1: 138:214695 REFERENCE 2: 138:214657 REFERENCE 3: 138:214651 REFERENCE 4: 138:214296 REFERENCE 5: 138:213944 REFERENCE 6: 138:212813 7: 138:212788 REFERENCE REFERENCE 8: 138:212782 REFERENCE 9: 138:209349 REFERENCE 10: 138:209218

=> d ide can 19

```
L9 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2003 ACS RN 109-89-7 REGISTRY CN Ethanamine, N-ethyl- (9CI) (CA INDEX NAME) OTHER CA INDEX NAMES:
```

sackey - 10 / 049463 CN Diethylamine (8CI) OTHER NAMES: CN DEA CN N, N-Diethylamine FS 3D CONCORD MF C4 H11 N CI COM LCSTN Files: AGRICOLA, ANABSTR, AQUIRE, BEILSTEIN\*, BIOBUSINESS, BIOSIS. BIOTECHNO, CA, CABA, CANCERLIT, CAOLD, CAPLUS, CASREACT, CBNB, CEN, CHEMCATS, CHEMINFORMRX, CHEMLIST, CHEMSAFE, CIN, CSCHEM, CSNB, DDFU, DETHERM\*, DIPPR\*, DRUGU, EMBASE, ENCOMPLIT, ENCOMPLIT2, ENCOMPPAT, ENCOMPPAT2, GMELIN\*, HODOC\*, HSDB\*, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE, MRCK\*, MSDS-OHS, NAPRALERT, NIOSHTIC, PDLCOM\*, PIRA, PROMT, RTECS\*, SPECINFO, SYNTHLINE, TOXCENTER, TULSA, ULIDAT, USPAT2, USPATFULL, VTB (\*File contains numerically searchable property data) Other Sources: DSL\*\*, EINECS\*\*, TSCA\*\* (\*\*Enter CHEMLIST File for up-to-date regulatory information) H3C-CH2-NH-CH2-CH3 \*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\* 14561 REFERENCES IN FILE CA (1962 TO DATE) 883 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA 14581 REFERENCES IN FILE CAPLUS (1962 TO DATE) 2 REFERENCES IN FILE CAOLD (PRIOR TO 1967) REFERENCE 1: 138:207281 REFERENCE 2: 138:205463 REFERENCE 3: 138:205229

REFERENCE 2: 138:205463

REFERENCE 3: 138:205229

REFERENCE 4: 138:205051

REFERENCE 5: 138:200070

REFERENCE 6: 138:197741

REFERENCE 7: 138:197696

REFERENCE 8: 138:197629

REFERENCE 9: 138:194052

REFERENCE 10: 138:192439

=> d ide can 110

L10 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2003 ACS RN 74-90-8 REGISTRY CN Hydrocyanic acid (8CI, 9CI) (CA INDEX NAME) OTHER NAMES: CN Carbon hydride nitride (CHN) CN Evercyn CN Formic anammonide CN Formonitrile CN Hydrogen cyanide

```
CN
     Prussic acid
CN
     Zyklon B
DR
     341972-31-4
MF
     CHN
CI
    COM
LC
                  ADISNEWS, AGRICOLA, ANABSTR, AQUIRE, BEILSTEIN*, BIOBUSINESS,
       BIOSIS, BIOTECHNO, CA, CABA, CANCERLIT, CAOLD, CAPLUS, CASREACT, CBNB,
       CEN, CHEMCATS, CHEMINFORMRX, CHEMLIST, CHEMSAFE, CIN, CSCHEM, CSNB,
       DDFU, DETHERM*, DIOGENES, DIPPR*, DRUGU, EMBASE, ENCOMPLIT, ENCOMPLIT2,
       ENCOMPPAT, ENCOMPPAT2, GMELIN*, HODOC*, HSDB*, IFICDB, IFIPAT, IFIUDB,
       IPA, MEDLINE, MRCK*, MSDS-OHS, NAPRALERT, NIOSHTIC, PDLCOM*, PIRA,
       PROMT, RTECS*, SPECINFO, TOXCENTER, TULSA, ULIDAT, USPAT2, USPATFULL,
         (*File contains numerically searchable property data)
                     DSL**, EINECS**, TSCA**
         (**Enter CHEMLIST File for up-to-date regulatory information)
 ,; N
HC
```

\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

1: 138:209211

11482 REFERENCES IN FILE CA (1962 TO DATE)
213 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
11489 REFERENCES IN FILE CAPLUS (1962 TO DATE)
2 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

REFERENCE 2: 138:206858 REFERENCE 3: 138:205462 REFERENCE 4: 138:201810 REFERENCE 138:195277 5: REFERENCE 6: 138:195216 REFERENCE 7: 138:195200 REFERENCE 8: 138:190465 REFERENCE 9: 138:186996 REFERENCE 10: 138:177624

REFERENCE

=> fil hcaplus FILE 'HCAPLUS' ENTERED AT 17:24:08 ON 02 APR 2003 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2003 AMERICAN CHEMICAL SOCIETY (ACS)

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FILE COVERS 1907 - 2 Apr 2003 VOL 138 ISS 14 FILE LAST UPDATED: 1 Apr 2003 (20030401/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

```
=> d all hitstr 122
L22 ANSWER 1 OF 1 HCAPLUS COPYRIGHT 2003 ACS
     2001:101098 HCAPLUS
ΑN
DN
     134:131253
ΤI
     Method and catalyst for making methyl ethyl
     ketone cyanohydrin
IN
     Croizy, Jean-francois; Esch, Marc; Esquirol, Gilbert
PΑ
     Atofina, Fr.
SO
     PCT Int. Appl., 13 pp.
     CODEN: PIXXD2
DT
     Patent
LA
     French
IC
     ICM C07C253-00
     ICS C07C255-12
     23-19 (Aliphatic Compounds)
     Section cross-reference(s): 45, 67
FAN.CNT 1
     PATENT NO.
                    KIND DATE
                                          APPLICATION NO. DATE
                                          -----
PΙ
     WO 2001009085
                     A1 20010208
                                         WO 2000-FR2136 20000725
        W: AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CU, CZ, DE,
             DK, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP,
             KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MD, MG, MK, MN,
             MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM,
             TR, TT, UA, UG, US, UZ, VN, YU, ZW, AM, AZ, BY, KG, KZ, MD, RU,
             TJ, TM
         RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY,
             DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ,
             CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG
     FR 2796939
                           20010202
                     A1
                                         FR 1999-9859
                                                           19990729
     FR 2796939
                      В1
                            20010914
     EP 1206445
                      A1
                            20020522
                                          EP 2000-958611
                                                           20000725
           AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT,
            IE, SI, LT, LV, FI, RO, MK, CY, AL
PRAI FR 1999-9859
                           19990729
                    . A
    WO 2000-FR2136
                     W
                           20000725
OS
    CASREACT 134:131253
AΒ
    Me Et ketone cyanohydrin is prepd. in high
     yield and selectivity by reacting hydrocyanic acid and
     2-butanone in the presence of catalytic amts. of diethylamine.
ST
    butanone cyanohydrin prepn; diethylamine hydrocyanation catalyst
    butanone cyanohydrin prepn
    Hydrocyanation catalysts
ΙT
        (diethylamine for the conversion of hydrocyanic
        acid and 2-butanone in the manuf. of Me Et
       ketone cyanohydrin)
ΙT
    Hydrocyanation
        (of hydrocyanic acid and 2-butanone in the manuf.
       of Me Et ketone cyanohydrin)
ΙT
    109-89-7, Diethylamine, uses
```

```
RL: CAT (Catalyst use); USES (Uses)
         (method and catalyst for making Me Et
         ketone cyanohydrin)
     74-90-8, Hydrogen cyanide, reactions
     78-93-3, Methyl ethyl ketone,
     reactions
     RL: RCT (Reactant); RACT (Reactant or reagent)
         (method and catalyst for making Me Et
         ketone cyanohydrin)
IT
     4111-08-4P, Methyl ethyl ketone
     cyanohydrin
     RL: SPN (Synthetic preparation); PREP (Preparation)
         (method and catalyst for making Me Et
        ketone cyanohydrin)
              THERE ARE 4 CITED REFERENCES AVAILABLE FOR THIS RECORD
RE.CNT
(1) Anon; HOUBEN-WEYL METHODEN DER ORGANISCHEN CHEMIE BAND VIII 1952, P274
(2) Saito, M; HCAPLUS
(3) Saito, M; JPN KOKAI TOKKYO KOHO P7
(4) Union Carbide Corp; WO 8500166 A 1985 HCAPLUS
     109-89-7, Diethylamine, uses
     RL: CAT (Catalyst use); USES (Uses)
         (method and catalyst for making {\tt Me} {\tt Et}
        ketone cyanohydrin)
RN
     109-89-7 HCAPLUS
     Ethanamine, N-ethyl- (9CI) (CA INDEX NAME)
CN
H3C-CH2-NH-CH2-CH3
ΙT
     74-90-8, Hydrogen cyanide, reactions
     78-93-3, Methyl ethyl ketone,
     reactions
     RL: RCT (Reactant); RACT (Reactant or reagent)
        (method and catalyst for making Me Et
        ketone cyanohydrin)
     74-90-8 HCAPLUS
RN
CN
     Hydrocyanic acid (8CI, 9CI) (CA INDEX NAME)
HC^{-\frac{1}{2}N}
RN
     78-93-3 HCAPLUS
CN
     2-Butanone (8CI, 9CI) (CA INDEX NAME)
H3C-C-CH2-CH3
     4111-08-4P, Methyl ethyl ketone
     cyanohydrin
     RL: SPN (Synthetic preparation); PREP (Preparation)
        (method and catalyst for making Me Et
        ketone cyanohydrin)
RN
     4111-08-4 HCAPLUS
CN
     Butanenitrile, 2-hydroxy-2-methyl- (9CI) (CA INDEX NAME)
```

#### => d bib abs hitstr tot 123

L23 ANSWER 1 OF 10 HCAPLUS COPYRIGHT 2003 ACS

AN 2001:224394 HCAPLUS

DN 134:233091

TI Preparation of plant-derived biopesticides and their synthetic analogs

IN Coats, Joel R.; Peterson, Christopher J.; Tsao, Rong; Eggler, Aimee L.; Tylka, Gregory L.

PA Iowa State University Research Foundation, Inc., USA

SO U.S., 18 pp. CODEN: USXXAM

DT Patent

LA English

FAN.CNT 1

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
US 6207705	В1	20010327	US 1997-828190	19970321
T 130 1000 10000	Б.	1000000		

PRAI US 1996-13956P P 19960322

AB Biopesticide compns. comprise a purified glucosinolate breakdown product wherein a starting material for the purified glucosinolate breakdown product is isolated from a crambe, rapeseed, flax, cassava, or mustard plant. Glucosinolate breakdown products, such as 1-cyano-2-hydroxy-3-butene, di-Me ketone cyanohydrin, Me Et ketone cyanohydrin, 3-cyano-3-hydroxy-1-propene, etc., and their analogs, as well as monoterpenoids are purified from plant exts. or synthetically prepd. The biopesticides are used against insects, mites, ticks, and nematodes as contact pesticides, aquatic pesticides, or fumigants.

IT 4111-08-4P, Methyl ethyl ketone cyanohydrin

RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); PUR (Purification or recovery); BIOL (Biological study); PREP (Preparation)

(prepn. of plant-derived biopesticides and their synthetic analogs)

RN 4111-08-4 HCAPLUS

CN Butanenitrile, 2-hydroxy-2-methyl- (9CI) (CA INDEX NAME)

## RE.CNT 53 THERE ARE 53 CITED REFERENCES AVAILABLE FOR THIS RECORD ALL CITATIONS AVAILABLE IN THE RE FORMAT

L23 ANSWER 2 OF 10 HCAPLUS COPYRIGHT 2003 ACS

AN 1996:154053 HCAPLUS

DN 124:224570

TI The first recombinant hydroxynitrile lyase and its application in the synthesis of (S)-cyanohydrins

AU Foerster, Siegfried; Roos, Juergen; Effenberger, Franz; Wajant, Harald; Sprauer, Achim

Inst. Org. Chemie Univ., Stuttgart, D-70569, Germany

Angewandte Chemie, International Edition in English (1996), 35(4), 437-9 CODEN: ACIEAY; ISSN: 0570-0833

PB VCH

DT Journal

LAEnglish

The authors overexpressed Manihot esculenta hydroxynitrile lyase (meHNL) AR in Escherichia coli. Enantioselective addn. of hydrocyanic acid to several aldehydes and ketones was by enzyme immobilized on nitrocellulose and using diisopropyl ether as solvent was demonstrated.

ΙT 174849-22-0P

RL: BPN (Biosynthetic preparation); BIOL (Biological study); PREP (Preparation)

(purifn. and characterization of recombinant hydroxynitrile lyase and its application in synthesis of (S)-cyanohydrins)

RN 174849-22-0 HCAPLUS

CN Butanenitrile, 2-hydroxy-2-methyl-, (2S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

ΙT **78-93-3,** 2-Butanone, reactions

RL: RCT (Reactant); RACT (Reactant or reagent) (purifn. and characterization of recombinant hydroxynitrile lyase and its application in synthesis of (S)-cyanohydrins)

RN 78-93-3 HCAPLUS

CN 2-Butanone (8CI, 9CI) (CA INDEX NAME)

L23 ANSWER 3 OF 10 HCAPLUS COPYRIGHT 2003 ACS

AN 1993:443947 HCAPLUS

DN 119:43947

TI Improved purification of an (R)-oxynitrilase from Linum usitatissimum (flax) and investigation of the substrate range

Albrecht, Jens; Jansen, Inge; Kula, Maria Regina AU

CS Inst. Enzymtechnol., Heinrich-Heine-Univ., Juelich, D-5170, Germany

SO Biotechnology and Applied Biochemistry (1993), 17(2), 191-203 CODEN: BABIEC; ISSN: 0885-4513

DTJournal

LA English

The purifn. of (R)-oxynitrilase (EC 4.1.2.10) from Linum usitatissimum has AB been improved considerably. The enzyme is obtained from seedlings in 60% yield by fractional salt pptn. followed by ion-exchange and hydrophobic-interaction chromatog. Final gel-permeation chromatog. yields a protein with a specific activity of 53 units/mg at pH 4.1. The N-terminal sequence is reported and microheterogeneity demonstrated. The substrate range was investigated using (R)-oxynitrilase immobilized on Eupergit and tert-Bu Me ether as solvent. The addn. of HCN to various aliph. ketones and aldehydes is catalyzed by the enzyme, while arom. ketones are not converted. (R)-butan-2-one cyanohydrin was synthesized on a preparative scale and the product characterized. ΙT

**122045-29-8P**, (R)-Butan-2-one cyanohydrin

```
RL: SPN (Synthetic preparation); PREP (Preparation)
                              (prepn. of, by immobilized oxynitrilase of flax)
  RN
                   122045-29-8 HCAPLUS
                   Butanenitrile, 2-hydroxy-2-methyl-, (2R)- (9CI) (CA INDEX NAME)
  CN
  Absolute stereochemistry.
    HO
                 CN
  Et R Me
                  74-90-8, Hydrogen cyanide, reactions
                  RL: RCT (Reactant); RACT (Reactant or reagent)
                             (reaction of, with aliph. ketones and aldehydes in oxynitrilase
                            presence)
                  74-90-8 HCAPLUS
 RN
 CN
                  Hydrocyanic acid (8CI, 9CI) (CA INDEX NAME)
                N
 HC '
                 78-93-3, Butanone, reactions
 IT
                 RL: RCT (Reactant); RACT (Reactant or reagent)
                            (reaction of, with oxynitrilase of Linum usitatissimum)
RN
                 78-93-3 HCAPLUS
CN
                 2-Butanone (8CI, 9CI) (CA INDEX NAME)
H3C-C-CH2-CH3
L23 ANSWER 4 OF 10 HCAPLUS COPYRIGHT 2003 ACS
                 1993:58234 HCAPLUS
AN
                 118:58234
DN
TI
                 Enzymic preparation of (R)-methyl and (R)-ethylketone cyanohydrins
ΙN
                 Effenberger, Franz; Ziegler, Thomas; Hoersch, Brigitte; Heid, Stephan
PA
                 Degussa A.-G., Germany
SO
                 Ger., 4 pp.
                 CODEN: GWXXAW
DT
                 Patent
                German
FAN.CNT 1
                 PATENT NO.
                 PATENT NO. KIND DATE
                                                                    KIND DATE
                                                                                                                                          APPLICATION NO. DATE
                                                                                                                                               -----
PΙ
                DE 4102327
                                                                        Cl 19920604
                                                                                                                                          DE 1991-4102327 19910126
PRAI DE 1991-4102327
                                                                                          19910126
               MARPAT 118:58234
                 (R)-Ketone cyanhydrins are prepd. for use in the manuf. of
                 .alpha.-hydroxy-.alpha.-ethyl- or .alpha.-hydroxy-.alpha.-methyl-
                carboxylic acids from the corresponding ketone and HCN using
                 (R)-oxynitrilase in an org. solvent. The enzyme is prepd. from almonds % \left( 1\right) =\left( 1\right) +\left( 
                and is preferably immobilized. Almond (R)-oxynitrilase 100 units (1000
                units/mL) was adsorbed onto Avicel cellulose 1 g (preswollen in Na acetate
                buffer pH 4.5) in diisopropylether 20 mL. Methylethyl
                ketone 5 mmol, and anhyd. HCN 400 .mu.L were added and
                the mixt. incubated at O.degree. for 4 h. Yield of (R)-
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methylethyl ketone cyanhydrin was 80% with an enantiomeric excess of 76%. 122045-29-8P IT RL: PREP (Preparation) (prepn. of, enzymic, from ketone, with immobilized oxynitrilase) 122045-29-8 HCAPLUS RN CN Butanenitrile, 2-hydroxy-2-methyl-, (2R)- (9CI) (CA INDEX NAME) Absolute stereochemistry. HO CN R Et Me IT 74-90-8, Hydrogen cyanide, reactions 78-93-3, Methylethyl ketone, reactions RL: RCT (Reactant); RACT (Reactant or reagent) (reactions of, in prepn. chiral cyanohydrins with immobilized oxynitrilase) 74-90-8 HCAPLUS RN Hydrocyanic acid (8CI, 9CI) (CA INDEX NAME) CN HC: N 78-93-3 HCAPLUS RN 2-Butanone (8CI, 9CI) (CA INDEX NAME) CN 0 H3C-C-CH2-CH3 L23 ANSWER 5 OF 10 HCAPLUS COPYRIGHT 2003 ACS 1991:491338 HCAPLUS AN DN 115:91338 Enzyme catalyzed reactions. 9. Enzyme-catalyzed synthesis of (R)-ketone TΙ cyanohydrins and their hydrolysis to (R)-.alpha.-hydroxy-.alpha.-methyl carboxylic acids Effenberger, Franz; Hoersch, Brigitte; Weingart, Franz; Ziegler, Thomas; ΑU Kuehner, Stefan Inst. Org. Chem., Univ. Stuttgart, Stuttgart, 7000/80, Germany CS Tetrahedron Letters (1991), 32(23), 2605-8 SO CODEN: TELEAY; ISSN: 0040-4039 DT Journal English LA CASREACT 115:91338 OS (R)-Ketone cyanohydrins (R)-HOCRMeCN (I) are obtained with high AB enantioselectivity from aliph. ketones and HCN in org. solvents using (R)-oxynitrilase (EC 4.1.2.10) as catalyst. Acid catalyzed hydrolysis of I affords the corresponding (R)-.alpha.-hydroxy-.alpha.methyl carboxylic acids without measurable racemization: ΙT 78-93-3, Ethyl methyl ketone, reactions RL: RCT (Reactant); RACT (Reactant or reagent) (enantioselective addn. of hydrogen cyanide to, in presence of oxynitrilase) 78-93-3 HCAPLUS RN

CN 2-Butanone (8CI, 9CI) (CA INDEX NAME)

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0
H3C-C-CH2-CH3
IT
     74-90-8, Hydrogen cyanide, reactions
     RL: RCT (Reactant); RACT (Reactant or reagent)
         (enantioselective addn. of, to ketones in presence of oxynitrilase)
RN
     74-90-8 HCAPLUS
     Hydrocyanic acid (8CI, 9CI) (CA INDEX NAME)
CN
HC 2
TΤ
     122045-29-8P
     RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT
     (Reactant or reagent)
         (prepn. and hydrolysis of)
RN
     122045-29-8 HCAPLUS
     Butanenitrile, 2-hydroxy-2-methyl-, (2R)- (9CI) (CA INDEX NAME)
CN
Absolute stereochemistry.
 HO CN
Et R Me
L23 ANSWER 6 OF 10 HCAPLUS COPYRIGHT 2003 ACS
AN
     1989:573178 HCAPLUS
DN
     111:173178
     Asymmetric hydrocyanation of a range of aromatic and aliphatic aldehydes
TI
     Matthews, Barry R.; Jackson, W. Roy; Jayatilake, Gamini S.; Wilshire,
ΑU
     Colin; Jacobs, Howard A.
CS
     Dep. Chem., Monash Univ., Clayton, 3168, Australia
SO
     Australian Journal of Chemistry (1988), 41(11), 1697-709
     CODEN: AJCHAS; ISSN: 0004-9425
DT
     Journal
LΑ
     English
OS
     CASREACT 111:173178
     A range of aryl, alkyl and heterocyclic aldehydes have been treated with
AB
     HCN in the presence of the Inoue catalyst, (R,R) - or
     (S,S)-cyclo[phenylalanylhistidyl]. Most aryl aldehydes with
     electron-donating substituents in the m- or p-positions give high
     enantiomeric excess (e.e.) values (.gtoreq.80%), but aryl aldehydes with
     strong electron-withdrawing substituents gave moderate e.e. values
     (.ltoreq.50%). These moderate values are believed to be due to partial
     racemization of the product cyanohydrins in the presence of the mildly
     basic catalyst. In contrast to the reactions of aryl aldehydes, reactions
     of alkyl aldehydes and of ketones gave low e.e. values (.ltoreq.30%), an
     explanation is proposed.
ΙT
     78-93-3, 2-Butanone, reactions
     RL: RCT (Reactant); RACT (Reactant or reagent)
        (hydrocyanation of, catalyst for)
RN
     78-93-3 HCAPLUS
     2-Butanone (8CI, 9CI) (CA INDEX NAME)
CN
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\cap
H3C-C-CH2-CH3
IT
     74-90-8
     RL: RCT (Reactant); RACT (Reactant or reagent)
        (hydrocyanation, stereoselective, of arom. and aliph. aldehydes)
     74-90-8 HCAPLUS
RN
     Hydrocyanic acid (8CI, 9CI) (CA INDEX NAME)
CN
HC<sup>2</sup>
IT
     122045-29-8P
     RL: SPN (Synthetic preparation); PREP (Preparation)
        (prepn. of)
RN
     122045-29-8 HCAPLUS
     Butanenitrile, 2-hydroxy-2-methyl-, (2R)- (9CI) (CA INDEX NAME)
CN
Absolute stereochemistry.
HO
     CN
Et R Me
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AN
     1985:199688 HCAPLUS
DN
     102:199688
     Properties of a microsomal enzyme system from Linum usitatissimum (linen
TΙ
     flax) which oxidizes valine to acetone cyanohydrin and isoleucine to
     2-methylbutanone cyanohydrin
     Cutler, Adrian J.; Sternberg, Margarete; Conn, Eric E. Dep. Biochem. Biophys., Univ. California, Davis, CA, 95616, USA
ΑU
CS
     Archives of Biochemistry and Biophysics (1985), 238(1), 272-9
SO
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ANSWER 7 OF 10 HCAPLUS COPYRIGHT 2003 ACS

CODEN: ABBIA4; ISSN: 0003-9861

DT Journal

LA English

AΒ Microsomal prepns. from flax seedlings have recently been shown to convert L-valine to acetone cyanohydrin, the precursor of the cyanogenic glucoside linamarin (Cutler, A. J.; Conn, E. E., 1981). Further details of this 4-step biosynthetic sequence and also details of the analogous reactions in lotaustralin biosynthesis have been obtained. The lotaustralin precursor, 2-methylbutyraldoxime, is the best substrate for cyanide prodn.  $(V_{\text{max}} = 413 \text{ nmol h-1 g fresh wt.-1})$  and inhibits the conversion of valine and isoleucine into products. Similarly, the linamarin precursor isobutyraldoxime is an excellent substrate (Vmax = 400 nmol h-1 g freshwt.-1) and also inhibits oxidn. of the amino acids. The substrate specificity of 'the oxime-metabolizing step is low and a variety of aliph. oximes are converted to cyanide. On the other hand, the activity of the microsomal ext. is highly selective with regard to the amino acid substrate since, of the aliph. amino acids tested, only valine and isoleucine are metabolized. Product formation from isobutyronitrile (a linamarin precursor) was not demonstrated, but detectable cyanide formation from 2-methylcyanobutane, the corresponding precursor of lotaustralin was obsd. Competition expts. showed that the biosynthesis of linamarin and lotaustralin is not likely to be catalyzed by sep. enzyme systems.

ΙT 4111-08-4P

> RL: BSU (Biological study, unclassified); MFM (Metabolic formation); BIOL (Biological study); FORM (Formation, nonpreparative); PREP (Preparation) (formation of, from isoleucine oxidn. by flax microsomes)

RN 4111-08-4 HCAPLUS

CN Butanenitrile, 2-hydroxy-2-methyl- (9CI) (CA INDEX NAME)

L23 ANSWER 8 OF 10 HCAPLUS COPYRIGHT 2003 ACS

1984:20417 HCAPLUS AN

100:20417 DN

ΤI Four aliphatic esters of Chamaemelum fuscatum essential oil

De Pascual-T., J.; Caballero, E.; Caballero, C.; Anaya, J.; Gonzalez, M. ΑU

CS

Dep. Org. Chem., Salamanca Univ., Salamanca, Spain Phytochemistry (Elsevier) (1983), 22(8), 1757-9 CODEN: PYTCAS; ISSN: 0031-9422 SO

DTJournal

LΑ English

AB In addn. to known compds., the esters H2C:CMeCO2CH2R [R = CMe:CHMe-(Z), CMe(OH)CH:CH2, CMe(OH)COMe] and neryl isovalerate were isolated from the essential oil of C. fuscatum and their structures were established by spectral methods and by synthesis.

ΙT 4111-08-4P

> RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(prepn. and hydrolysis-elimination reaction of)

4111-08-4 HCAPLUS RN

CNButanenitrile, 2-hydroxy-2-methyl- (9CI)

IT **78-93-3**, reactions

RL: RCT (Reactant); RACT (Reactant or reagent) (reaction of, with sodium cyanide)

78-93-3 HCAPLUS RN

CN 2-Butanone (8CI, 9CI) (CA INDEX NAME)

L23 ANSWER 9 OF 10 HCAPLUS COPYRIGHT 2003 ACS AN 1982:6417 HCAPLUS

DN 96:6417

TI Cyanomethyl .alpha.-(p-phenoxyphenoxy)propionates

PA Compagnie Francaise de Produits Industriels, Fr.

SO Fr. Demande, 41 pp. Addn. to Fr. Appl. No. 79 01,020. CODEN: FRXXBL

DT Patent

LA French

FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI PRAI GI	FR 2473514 US 1980-112352	A2	19810717 19800115	FR 1981-666	19810115

$$\begin{array}{c|c} & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & \\ & & & \\ & &$$

AB Title esters I [n = 1, 2; R = Cl, CF3; R1 and R2 (same or different) are H, alkyl, chlorophenyl, (trifluoromethyl)phenyl, or R1R2 = polymethylene], useful as herbicides, were prepd. by different methods. Thus, 4-(4-F3CC6H4O)C6H4OCHMeCO2H was converted to its acid chloride, which was esterified with Me2C(OH)CN and pyridine to give I (Rn = 4-F3C, R1 = R2 = Me). Herbicidal activity data are presented for I.

Ι

IT **78-93-3**, reactions

RL: RCT (Reactant); RACT (Reactant or reagent)
 (conversion of, to cyanohydrin)

RN 78-93-3 HCAPLUS

CN 2-Butanone (8CI, 9CI) (CA INDEX NAME)

#### IT 4111-08-4P

RL: SPN (Synthetic preparation); PREP (Preparation) (prepn. of, and esterification of .alpha.-(p-phenoxyphenoxy)propionyl chloride deriv. by)

RN 4111-08-4 HCAPLUS

CN Butanenitrile, 2-hydroxy-2-methyl- (9CI) (CA INDEX NAME)

L23 · ANSWER 10 OF 10 HCAPLUS COPYRIGHT 2003 ACS

AN 1972:58924 HCAPLUS

DN 76:58924

TI Synthesis of unsaturated aliphatic nitriles

AU Mekhtiev, S. I.; Mamedov, R. G.

CS VNIIolefin, Baku, USSR

SO Azerbaidzhanskii Khimicheskii Zhurnal (1971), (2), 110-15

CODEN: AZKZAU; ISSN: 0005-2531

DT Journal

LA Russian

AB Cyanohydrins (I) (80-95% based on HCN) were obtained by hydrocyanidation of ketones in the presence of KOH. The stability of I in alk. medium increased with increase of their mol. wt. Alkylacrylonitriles (II) were prepd. by dehydrating I from MeCOEt, Et2CO, MeCOPr, and MeCOBu. II (80-2% based on converted I) were obtained at I-P2O5-quinoline mole ratio 1:1.25:1, 60-80.degree., and reaction time 1 hr.

IT 4111-08-4P

RN 4111-08-4 HCAPLUS

CN Butanenitrile, 2-hydroxy-2-methyl- (9CI) (CA INDEX NAME)

IT 74-90-8

RL: RCT (Reactant); RACT (Reactant or reagent)
 (reaction of, with aliphatic ketones)

RN 74-90-8 HCAPLUS

CN Hydrocyanic acid (8CI, 9CI) (CA INDEX NAME)

IT **78-93-3**, reactions

RL: RCT (Reactant); RACT (Reactant or reagent)
 (with hydrocyanic acid)

RN 78-93-3 HCAPLUS

CN 2-Butanone (8CI, 9CI) (CA INDEX NAME)